

10/549,852 Yong Chu X=S\_alkyl 05/15/2009 Yong Chu

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NEWS 16 MAR 11 EPFULL backfile enhanced with additional full-text  
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NEWS 17 MAR 11 ESBIOBASE reloaded and enhanced  
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NEWS 19 MAR 23 CA/CAPLUS enhanced with more than 250,000 patent  
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NEWS 21 APR 03 CAS coverage of exemplified prophetic substances  
enhanced  
NEWS 22 APR 07 STN is raising the limits on saved answers  
NEWS 23 APR 24 CA/CAPLUS now has more comprehensive patent assignee  
information  
NEWS 24 APR 26 USPATFULL and USPAT2 enhanced with patent  
assignment/reassignment information  
NEWS 25 APR 28 CAS patent authority coverage expanded  
NEWS 26 APR 28 ENCOMPLIT/ENCOMPLIT2 search fields enhanced  
NEWS 27 APR 28 Limits doubled for structure searching in CAS

REGISTRY

NEWS 28 MAY 08 STN Express, Version 8.4, now available

NEWS 29 MAY 11 STN on the Web enhanced

NEWS 30 MAY 11 BEILSTEIN substance information now available on STN Easy

NEWS 31 MAY 14 DGENE, PCTGEN and USGENE enhanced with increased limits for exact sequence match searches and introduction of free HIT display format

NEWS 32 MAY 15 INPADOCDB and INPAFAMDB enhanced with Chinese legal status data

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,  
AND CURRENT DISCOVER FILE IS DATED 06 APRIL 2009.

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\* \* \* \* \* STN Columbus \* \* \* \* \*

FILE 'HOME' ENTERED AT 16:15:25 ON 15 MAY 2009

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	ENTRY	SESSION
FULL ESTIMATED COST	0.22	0.22

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STRUCTURE FILE UPDATES: 14 MAY 2009 HIGHEST RN 1146852-72-3  
DICTIONARY FILE UPDATES: 14 MAY 2009 HIGHEST RN 1146852-72-3

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TSCA INFORMATION NOW CURRENT THROUGH January 9, 2009.

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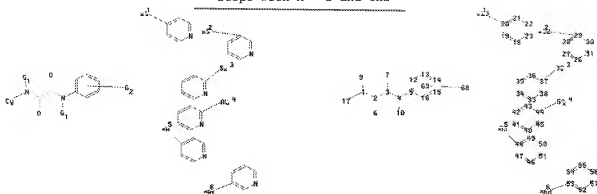
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\*\*Scope with X = S and CH2\*\*



chain nodes :

1 2 3 4 6 7 9 10 17 25 32 39 58 59 60 68

ring nodes :

5 12 13 14 15 16 18 19 20 21 22 23 26 27 28 29 30 31 33 34 35  
36 37 38 40 41 42 43 44 45 46 47 48 49 50 51 52 53 54 55 56 57

chain bonds :

1-9 1-2 1-17 2-3 2-6 3-4 3-7 4-5 4-10 20-25 29-32 37-39 44-58 48-59  
53-60

ring bonds :

5-12 5-16 12-13 13-14 14-15 15-16 18-19 18-23 19-20 20-21 21-22 22-23  
26-27 26-31 27-28 28-29 29-30 30-31 33-34 33-38 34-35 35-36 36-37 37-38  
40-41 40-45  
41-42 42-43 43-44 44-45 46-47 46-51 47-48 48-49 49-50 50-51 52-53 52-57  
53-54 54-55 55-56  
56-57

exact/norm bonds :

1-9 1-2 1-17 2-3 2-6 3-4 3-7 4-5 4-10 20-25 29-32 37-39 44-58 48-59  
53-60

normalized bonds :

5-12 5-16 12-13 13-14 14-15 15-16 18-19 18-23 19-20 20-21 21-22 22-23  
26-27 26-31 27-28 28-29 29-30 30-31 33-34 33-38 34-35 35-36 36-37 37-38  
40-41 40-45  
41-42 42-43 43-44 44-45 46-47 46-51 47-48 48-49 49-50 50-51 52-53 52-57  
53-54 54-55 55-56  
56-57

G1:H,CH3,CH2,Et

G2:[\*1],[\*2],[\*3],[\*4],[\*5],[\*6]

Connectivity :

6:1 E exact RC ring/chain 7:1 E exact RC ring/chain

Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:Atom 6:CLASS 7:CLASS 9:CLASS 10:CLASS  
12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom  
21:Atom 22:Atom  
23:Atom 25:CLASS 26:Atom 27:Atom 28:Atom 29:Atom 30:Atom 31:Atom 32:CLASS  
33:Atom 34:Atom

35:Atom 36:Atom 37:Atom 38:Atom 39:CLASS 40:Atom 41:Atom 42:Atom 43:Atom  
 44:Atom 45:Atom  
 46:Atom 47:Atom 48:Atom 49:Atom 50:Atom 51:Atom 52:Atom 53:Atom 54:Atom  
 55:Atom 56:Atom  
 57:Atom 58:CLASS 59:CLASS 60:CLASS 68:CLASS 69:Atom  
 Generic attributes :  
 17:  
 Saturation : Unsaturated  
 58:  
 Saturation : Saturated  
 59:  
 Saturation : Saturated  
 60:  
 Saturation : Saturated

L1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1 STR

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

Structure attributes must be viewed using STN Express query preparation.

=> s l1

SAMPLE SEARCH INITIATED 16:16:27 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 508 TO ITERATE

100.0% PROCESSED 508 ITERATIONS

0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*

BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 8808 TO 11512

PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 full

FULL SEARCH INITIATED 16:16:40 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 9677 TO ITERATE

100.0% PROCESSED 9677 ITERATIONS

5 ANSWERS

SEARCH TIME: 00.00.01

L3 5 SEA SSS FUL L1

=> file caplus

COST IN U.S. DOLLARS

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TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

186.36

186.58

FILE 'CAPLUS' ENTERED AT 16:16:48 ON 15 MAY 2009

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FILE COVERS 1907 - 15 May 2009 VOL 150 ISS 21  
FILE LAST UPDATED: 14 May 2009 (20090514/ED)  
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Feb 2009  
USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Feb 2009

CAPLUS now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

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This file contains CAS Registry Numbers for easy and accurate

=> s l3

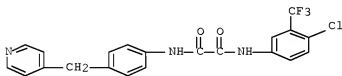
L4 2 L3

=> d ibib abs hitstr tot

L4 ANSWER 1 OF 2 CAPLUS COPYRIGHT 2009 ACS on STN  
ACCESSION NUMBER: 2004:817864 CAPLUS Full-text  
DOCUMENT NUMBER: 141:314164  
TITLE: Preparation of pyridinyloxyphenylethanedi-  
amides as RAF-kinase inhibitors  
INVENTOR(S): Buchstaller, Hans-Peter; Wiesner, Matthias; Zenke,  
Frank; Amendt, Christiane; Grell, Matthias;  
Sirrenberg, Christian  
PATENT ASSIGNEE(S): Merck Patent G.m.b.H., Germany  
SOURCE: PCT Int. Appl., 197 pp.  
CODEN: PIXXD2  
DOCUMENT TYPE: Patent  
LANGUAGE: English  
FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004085399	A1	20041007	WO 2004-EP2406	20040309
W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, GR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
RW:	BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE,			

ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GO, GW, ML, MR, NE, SN, TD, TG			
AU 2004224239	A1	20041007	AU 2004-224239
CA 2520009	A1	20041007	CA 2004-2520009
EP 1606260	A1	20051221	EP 2004-718645
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK			
BR 2004007968	A	20060307	BR 2004-7968
CN 1764645	A	20060426	CN 2004-80007867
JP 2006521304	T	20060921	JP 2006-504685
US 20060189665	A1	20060824	<u>US 2005-549852</u>
ZA 2005008522	A	20070425	ZA 2005-8522
PRIORITY APPLN. INFO.:			EP 2003-6702
			WO 2004-EP2406
OTHER SOURCE(S):			CASREACT 141:314164; MARPAT 141:314164
AB	ADB [D = (substituted) bivalent oxamide moiety; A = L(ML)a; L = 5-7 membered cyclic structure, preferably aryl, heteroaryl, arylene, heteroarylene; Ll = (substituted) cyclic moiety having at least 5 members, preferably aryl, heteroaryl, aralkyl, cycloalkyl, heterocyclyl; M = bond, bridging group; a = 1-4; L, Ll contain 0-4 N, O, S atoms; B = (substituted) up to tricyclic aryl, heteroaryl contg. 0-4 N, O, S atoms], were prepd. for treatment of hyperproliferative and nonhyperproliferative disorders (no data). For example, reaction of N-(4-chloro-3-trifluoromethylphenyl)-2-oxoglycine (prepn. given) with 4-(4-pyridinyloxy)phenylamine yielded N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-(4-(4-pyridinyloxy)phenyl)ethanediamine.		
IT	767358-38-3P RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (prepn. of pyridinyloxyphenylethanediamide derivs. as RAF-kinase inhibitors)		
RN	767358-38-3 CAPLUS		
CN	Ethanediameide, N1-[4-chloro-3-(trifluoromethyl)phenyl]-N2-[4-(4-pyridinylmethyl)phenyl]- (CA INDEX NAME)		



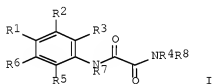
REFERENCE COUNT: 7 THERE ARE 7 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2009 ACS on SIN  
 ACCESSION NUMBER: 2001:631913 CAPLUS Full-text  
 DOCUMENT NUMBER: 135:195556  
 TITLE: Preparation of azolyphenyl oxamides as inosine monophosphate dehydrogenase (IMPDH) inhibitors  
 INVENTOR(S): Broadhurst, Michael John; Hill, Christopher Huw; Hurst, David Nigel; Jones, Philip Stephen; Kay, Paul Brittain; Kilford, Ian Reginald; Mckinnell, Robert Murray  
 PATENT ASSIGNEE(S): F. Hoffmann-La Roche A.-G., Switz.  
 SOURCE: Eur. Pat. Appl., 256 pp.

DOCUMENT TYPE: CODEN: EPXXDW  
 LANGUAGE: Patent  
 FAMILY ACC. NUM. COUNT: English  
 PATENT INFORMATION: 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
EP 1127883	A2	20010829	EP 2001-103521	20010216
EP 1127883	A3	20020807		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO				
US 20020052513	A1	20020502	US 2001-779116	20010208
US 6867299	B2	20050315		
CA 2337588	A1	20010824	CA 2001-2337588	20010220
HU 2001000836	A2	20011028	HU 2001-836	20010221
HR 2001000127	A1	20011231	HR 2001-127	20010221
NO 2001000900	A	20010827	NO 2001-900	20010222
CN 1310179	A	20010829	CN 2001-104906	20010223
BR 2001000790	A	20010925	BR 2001-790	20010223
IN 2001MA00167	A	20050304	IN 2001-MA167	20010223
JP 2001261663	A	20010926	JP 2001-51064	20010226
PRIORITY APPLN. INFO.:			GB 2000-4392	A 20000224
			GB 2000-15877	A 20000628
			GB 2000-20322	A 20000817

OTHER SOURCE(S): MARPAT 135:195556  
 GI



AB Title compds. (I; R1 = heterocyclyl; R2 = H, alkyl, alkoxy, halo, OH, cyano; R3 = H, alkyl, alkoxy, halo, cyano; R4 = H, alkyl, cycloalkyl, aryl, heterocyclyl; R5 = H, alkyl, alkoxy, halo, cyano; R6 = H, alkyl, alkoxy, halo, cyano; R7, R8 = H, alkyl; R4R8N = heterocyclyl), were prepd. Thus, 1,1-dimethyl-3-(4-nitrophenoxy)propylamine (prepn. given) was coupled with N-[3-methoxy-4-(5-oxazolyl)phenyl]oxamic acid in the presence of 1-(3-dimethylaminopropyl)-3-ethylcarbodiimide and 1-hydroxy-7-azabenzotriazole to give N-[3-methoxy-4-(5-oxazolyl)phenyl]-N'-[1,1-dimethyl-3-(4-nitrophenoxy)propyl]oxalamide. Tested I inhibited IMPDH with IC50 = 0.010-0.277 .mu.M. I can be used for treating immune mediated conditions or diseases, viral diseases, bacterial diseases, parasitic diseases, inflammation, inflammatory diseases, hyperproliferative vascular diseases, tumors, and cancer.

IT 357184-58-8P  
 RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

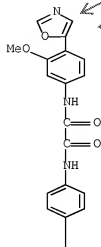
(prepn. of azolylphenyl oxamides as inosine monophosphate dehydrogenase (IMPDH) inhibitors)

RN 357184-58-8 CAPLUS

CN Ethanediamide, N1-[3-methoxy-4-(5-oxazolyl)phenyl]-N2-[4-(4-pyridinylmethyl)phenyl]- (CA INDEX NAME)

Close but not a prior art

\*\*Different R<sub>8</sub> PAGE 1-A



PAGE 2-A



REFERENCE COUNT: 9 THERE ARE 9 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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=>

Executing the logoff script...

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COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

12.28

198.86

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE

TOTAL



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ENTRY	SESSION
-1.64	-1.64

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